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|----------------------------|---------------|---------------|------------------|
| Species Tag: | 40002 | Species Name: | NaOH |
| Version: | 1 | | Sodium hydroxide |
| Date: | Aug. 1987 | | |
| Contributor: | R. L. Poynter | | |
| Lines Listed: | 91 | Q(300.0)= | 498.884 |
| Freq. (GHz) < | 2201 | Q(225.0)= | 374.025 |
| Max. J: | 91 | Q(150.0)= | 249.345 |
| LOGSTR0= | -6.3 | Q(75.00)= | 124.767 |
| LOGSTR1= | -8.0 | Q(37.50)= | 62.532 |
| Isotope Corr.: | 0.0 | Q(18.75)= | 31.427 |
| Egy. (cm ⁻¹) > | 0.0 | Q(9.375)= | 15.882 |
| μ_a = | 1.0 | A= | |
| μ_b = | | B= | 12567.0590(120) |
| μ_c = | | C= | |

The measured lines were taken from E. F. Pearson and M. B. Trueblood, 1973, *Astrophys. J. Lett. Ed.* **179**, L146, and P. Kuipers *et al.*, 1976, *Chem. Phys.* **15**, 457. Because no dipole moment measurement has been reported, we assumed a value of unity in our calculation. Two theoretical estimates have been published: N. G. Rambidi, Yu. G. Abashkin, and A. I. Dement'ev, 1984, *Russ. J. Inorg. Chem.* **29**, 12-22, and G. A. Long, J. F. Capitani, and L. Pedersen, 1983, *J. Mol. Struct.* **105**, 229-230. Of these, the Rambidi *et al.* value is closer to our estimate of 6.63 Debye with an estimated error of maybe 1 Debye, which is based upon extrapolations from the dipole moments of Li, Cs, and Na fluorides and the Cs and Li hydroxides. The Rambidi *et al.* value is 6.706 debye, while the Long *et al.* value ranges between 6.30 and 5.42 debye, depending on the basis set used in their calculations. The line intensities should be multiplied by a factor of about 44 to get something closer to the right value. The theoretical calculations suggest that this molecule is linear. A measurement is needed.